

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061119\
 Data File : BN006255.D
 Acq On : 11 Jun 2019 20:22
 Operator : JU/SJ
 Sample : K3083-05
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A51F8

Manual Integrations
 APPROVED

mohammad
 6/12/2019 8:19:07 AM

Quant Time: Jun 12 02:26:48 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jun 09 05:10:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	6312	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	25085	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	12773	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	23477m	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	14008m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	21526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	7177	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	11883m	0.19	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.01	152	6847	0.096	ng/ul	96
8) Acenaphthene	14.36	153	2410	0.049	ng/ul	98
9) Fluorene	15.35	166	3985	0.069	ng/ul#	95
12) Phenanthrene	17.09	178	192160m	2.550	ng/ul	
13) Anthracene	17.18	178	15826m	0.220	ng/ul	
15) Fluoranthene	19.12	202	662596m	7.962	ng/ul	
17) Pyrene	19.49	202	496902	6.734	ng/ul	100
18) Benzo(a)anthracene	21.25	228	181427m	2.928	ng/ul	
19) Chrysene	21.30	228	282493	4.857	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	396969m	4.423	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	134671m	1.540	ng/ul	
23) Benzo(a)pyrene	23.40	252	226816	2.713	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.73	276	204298	2.258	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.73	278	43272	0.598	ng/ul	98
26) Benzo(g,h,i)perylene	26.41	276	174156	2.298	ng/ul	94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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